

Spin-Wave Relaxation in a Quantum Hall Ferromagnet

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(Dated: December 9, 2008)

We study spin wave relaxation in quantum Hall ferromagnet regimes. Spin-orbit coupling is considered as a factor determining spin nonconservation, and external random potential as a cause of energy dissipation making spin-flip processes irreversible. We compare this relaxation mechanism with other relaxation channels existing in a quantum Hall ferromagnet.

PACS numbers 73.21.Fg, 73.43.Lp, 78.67.De

1. Last years are characterized by growing interest in spin relaxation (SR) in low-dimension systems — first of all, in the relaxation in quantum dots studied within the projects aimed at development of a computer employing spin memory. Yet, the relaxation of an electron spin in lateral quantum dots manufactured on the basis of two-dimensional (2D) heterostructures, should be in many respects similar to the SR of electrons localized in the 2D layer in minima of a smooth random potential (SRP). In high magnetic fields this single-electron relaxation corresponds to the situation occurring at low Landau level (LL) filling: $\nu \ll 1$ or $|\nu - 2n| \ll 1$ (n is an integer).¹

The SR at different filling factors, $\nu \gtrsim 1$, has quite different nature representing in this case a many-electron process. In particular, in a quantum Hall ferromagnet (QHF), i.e. at $\nu = 1, 3, \dots$ or $\nu = 1/3, 1/5, \dots$, the SR reduces to the relaxation of lowest collective excitations, i.e. spin waves.^{2,3} The SR observation would thereby be a good tool to study fundamental collective properties of a strongly correlated 2D electron gas (2DEG). However, in spite of much recent interest in the SR in a 2DEG, up to now only a handful of experiments relevant to the SR in a QHF were performed: these are indirect results based on the linewidth measurements in the electron spin resonance,⁴ and a direct observation where the photoluminescence dynamics of spin-up and spin-down states was studied.⁵ Meanwhile, availability of the new time-resolved technique of photon counting allows us to believe that new direct experiments on observation of excitations' relaxation in a 2DEG, in particular of the spin wave relaxation (SWR), will become available in the near future.⁶

Theoretically the SWR in a QHF was studied in works 7,8. It is worth noting here that

the SWR represents actually not spin dephasing but the energy relaxation due to the spin-flip process. Indeed, any spin-flip means at least dissipation of the Zeeman energy $\epsilon_Z = |g|\mu_B B$ ($g \approx -0.44$ in a GaAs structure). The latter is a part of the spin-wave (spin exciton, SE) energy

$$E_{\text{sw}} = \epsilon_Z + \mathcal{E}_q, \quad (1)$$

where \mathcal{E}_q is the SE correlation energy depending on the 2D wave vector q .^{2,3} At variance with the relaxation channel of Ref. 7 where electron-phonon interaction was considered as the mechanism making the relaxation irreversible, and contrary to the case of Ref. 8 where the irreversibility was provided by an inter-spin-exciton interaction mechanism, we now study smooth disorder field as the reason causing the energy transform. The SRP thereby determines an alternative relaxation channel competing with the ones studied earlier. Another distinction of the present work from Refs. 7,8 consists in the study of not only the integer QHF (at $\nu = 1, 3, \dots$) but also of the fractional one ($\nu = 1/3, 1/5, \dots$) as well. At the same time we again consider the spin-orbit coupling (SO) as the cause mixing different spin states and therefore providing the spin nonconservation. Actually, various SWR channels coexist in parallel. We consider the total rate and find crossover regions of external parameters (magnetic field, temperature, etc.) where one relaxation channel ceases to be dominant and changes into another.

The SR channel due to SRP was already considered in the integer quantum Hall ferromagnetic case.^{1,9} However, studied in these works instead of the SWR was a specific SR when initially the total macroscopic spin \vec{S} of the system as a whole is turned away from the equilibrium direction parallel to \vec{B} . (Relaxation of this Goldstone mode microscopically reduces to annihilation processes of the so-called zero SEs, having exactly zero momenta.) Contrary to this case, the spin perturbation determined by excitation of the spin waves (non-zero SEs) represents an initial deviation where $\Delta S = \Delta S_z$, so that \vec{S} is kept parallel to \vec{B} and the total symmetry of system remains unchanged.

Concerning the origin of SRP, one should note that it has in the 2D layer the “direct” component and the effective one. The former is the SRP determined by charged donors located outside the spacer. The latter is essential in some kinds of quantum wells, being determined by spatial fluctuations (in the plane of the layer) of quantum well width. These fluctuations lead to fluctuations of the size-quantization energy and may be presented as an SRP term in the single electron Hamiltonian. Both SRP components have approximately the same amplitude $\Delta \sim 10$ K and correlation length $\Lambda \sim 30 - 50$ nm.

2. The total Hamiltonian has form $H_{\text{tot}} = \sum_j H_1^{(j)} + H_{\text{int}}$, where j enumerates electrons,

H_{int} is the e - e interaction, and the single-electron operator is

$$H_1 = \hbar^2 \hat{\mathbf{q}}^2 / 2m_e^* - \epsilon_Z \hat{\sigma}_z / 2 + H_{SO} + \varphi(\mathbf{r}). \quad (2)$$

In this equation $\varphi(\mathbf{r})$ is the SRP field; the SO Hamiltonian is specified for the (001) GaAs plane,

$$H_{SO} = \alpha (\hat{\mathbf{q}} \times \hat{\boldsymbol{\sigma}})_z + \beta (\hat{q}_y \hat{\sigma}_y - \hat{q}_x \hat{\sigma}_x), \quad (3)$$

presenting a combination of the Rashba term and the crystalline anisotropy term¹⁰ ($\hat{\mathbf{q}} = -i\nabla + e\mathbf{A}/c\hbar$ is a 2D operator, $\sigma_{x,y,z}$ are the Pauli matrices). If the SRP is assumed to be Gaussian, then it is defined by the correlator $K(\mathbf{r}) = \langle \varphi(\mathbf{r}) \varphi(0) \rangle$. By choosing $\langle \varphi(\mathbf{r}) \rangle = 0$, in terms of the correlation length Λ and the LL width Δ the correlator is

$$K(\mathbf{r}) = \Delta^2 \exp(-r^2/\Lambda^2). \quad (4)$$

We first find the bare single-electron basis diagonalizing the Hamiltonian (2) without the SRP field. To within the leading order in the H_{SO} terms we obtain

$$\begin{aligned} \Psi_{pa} &= \begin{pmatrix} \psi_{np} \\ v\sqrt{n+1}\psi_{n+1,p} + iu\sqrt{n}\psi_{n-1,p} \end{pmatrix}, \\ \Psi_{pb} &= \begin{pmatrix} -v\sqrt{n}\psi_{n-1,p} + iu\sqrt{n+1}\psi_{n+1,p} \\ \psi_{np} \end{pmatrix} \end{aligned} \quad (5)$$

Here ψ_{np} is the electron wave function in the Landau gauge, n is the number of the half-filled LL in the odd-integer quantum Hall regime, i.e. in the $\nu = 2n+1$ case. Otherwise, if $\nu \leq 1$, we set $n=0$. u and v are small dimensionless parameters: $u = \beta\sqrt{2}/l_B\hbar\omega_c$ and $v = \alpha\sqrt{2}/l_B\hbar\omega_c$ (ω_c and l_B are the cyclotron frequency and the magnetic length, respectively). The single-electron states thus cease to be purely spin states but acquire a chirality a or b . The spin flip corresponds thereby to the $a \rightarrow b$ process now.

By analogy with previous works^{1,7,8,9} (see also Ref. 11) we define the SE creation operator

$$\mathcal{Q}_{ab\mathbf{q}}^\dagger = \frac{1}{\sqrt{N_\phi}} \sum_p e^{-iq_x p} b_{p+\frac{q_y}{2}}^\dagger a_{p-\frac{q_y}{2}}, \quad (6)$$

where a_p and b_p are the Fermi annihilation operators corresponding to states (5), N_ϕ is the LL degeneracy number. In Eq. (6) and everywhere below we measure wave vector q in the $1/l_B$ units. If the ratio $r_c = (\alpha e^2/\kappa l_B)/\hbar\omega_c$ is considered to be small ($\alpha < 1$ is the averaged formfactor which appears due to finiteness of the layer thickness), and the SRP and SO terms in Eq. (2) are ignored, then the operator (6) acting on the ground state in the *odd-integer quantum Hall regime* yields the *eigen state* of the total Hamiltonian: namely,

$[H_{\text{tot}}, \mathcal{Q}_{ab\mathbf{q}}^\dagger]|0\rangle = (\epsilon_Z + \mathcal{E}_q)\mathcal{Q}_{ab\mathbf{q}}^\dagger|0\rangle$, where $|0\rangle = |\underbrace{\uparrow, \uparrow, \dots \uparrow}_{N_\phi}\rangle$. This basic property of the exciton state, $\mathcal{Q}_{ab\mathbf{q}}^\dagger|0\rangle$, is the asymptotically exact one to the first order in r_c .

Now consider corrections arising due to the H_{SO} terms. When presented in terms of basis states (5), spin operators $\int \Psi^\dagger \hat{\mathbf{S}}^2 \Psi d^2\mathbf{r}$ and $\int \Psi^\dagger \hat{S}_z \Psi d^2\mathbf{r}$ [where $\Psi = \sum_p (a_p \Psi_{pa} + b_p \Psi_{pb})$] preserve invariant form up to the second order in u and v . However, the interaction Hamiltonian $H_{\text{int}} = \frac{1}{2} \int d\mathbf{r}_1 d\mathbf{r}_2 \Psi^\dagger(\mathbf{r}_2) \Psi^\dagger(\mathbf{r}_1) U(\mathbf{r}_1 - \mathbf{r}_2) \Psi(\mathbf{r}_1) \Psi(\mathbf{r}_2)$ acquires proportional to u and v terms which correspond to creation and annihilation of SEs in the system. It is exactly these terms that lead to the “coalescence” channel of the SWR.⁸ In the present work we study another relaxation channel. Therefore, neglecting this SO corrections to \hat{H}_{int} , we focus on the SRP term. Calculating $\int \Psi^\dagger \varphi(\mathbf{r}) \Psi d^2\mathbf{r}$, we get the terms responsible for a spin-flip:

$$\hat{\varphi} = N_\phi^{1/2} l_B \sum_{\mathbf{q}} \bar{\varphi}(\mathbf{q}) (iuq_+ - vq_-) \mathcal{Q}_{\mathbf{q}} + \text{H.c.} \quad (7)$$

(it is assumed here that $q \ll 1$). $\bar{\varphi}(\mathbf{q})$ is the Fourier component [i.e. $\varphi = \sum_{\mathbf{q}} \bar{\varphi}(\mathbf{q}) e^{i\mathbf{q}\mathbf{r}}$], and $q_{\pm} = \mp i(q_x \pm iq_y)/\sqrt{2}$.

At variance with integer QHF, the use of the excitonic basis $\mathcal{Q}_{ab\mathbf{q}}^\dagger |0\rangle$ presents only a *model approach* in the case of *fractional quantum Hall regime*. Generally, spin-flip excitations within the same Landau level might be many-particle rather than two-particle excitations at fractional filling because the same change of the spin numbers $\delta S = \delta S_z = -1$ may be achieved with participation of arbitrary number of intra-spin-sublevel excitations (charge-density waves). These waves are generated by the operator $\mathcal{A}_{\mathbf{q}}^\dagger = N_\phi^{-1/2} \mathcal{Q}_{aa\mathbf{q}}^\dagger$ acting on the ground state $|0\rangle = |\overbrace{\uparrow, \dots, \uparrow}^{\nu N_\phi}, \dots, \uparrow\rangle$.¹² It is trivial in the case of integer ν ($\mathcal{A}_{\mathbf{q}}^\dagger |0\rangle = \delta_{\mathbf{q},0} |0\rangle$); however, states of the $\mathcal{Q}_{ab\mathbf{q}_1}^\dagger \mathcal{A}_{\mathbf{q}_2}^\dagger \mathcal{A}_{\mathbf{q}_3}^\dagger \dots |0\rangle$ type might constitute a basis set if one studies a spin-flip at fractional ν . On the other hand, a comprehensive phenomenological analysis^{3,12} suggests that even the spin-flip basis reduced to single-mode (single-exciton) states would be quite appropriate, at least for lowest-energy excitations in the case of fractional QHF. This single-mode approach is indirectly substantiated by the fact that the charge-density wave has a Coulomb gap¹² which is well larger than the Zeeman gap ϵ_Z . Hence for a fractional QHF, just as in Ref. 3, we will consider the only state $\mathcal{Q}_{ab\mathbf{q}}^\dagger |0\rangle$ to describe the spin-flip excitation. The commutation algebra for operators $\mathcal{Q}_{ab\mathbf{q}}^\dagger$, $\mathcal{A}_{\mathbf{q}'}^\dagger$ and $\mathcal{B}_{\mathbf{q}''}^\dagger = N_\phi^{-1/2} \mathcal{Q}_{bb\mathbf{q}''}^\dagger$ is certainly the same as for integer filling,^{7,8,9}. However, a difference arises in the calculation of expectation $\langle 0 | \mathcal{A}_{\mathbf{q}} \mathcal{A}_{\mathbf{q}'}^\dagger | 0 \rangle$ which is needful for the following. This value is simply $\delta_{\mathbf{q},0} \delta_{\mathbf{q}',0}$ at integer filling, but at $\nu < 1$ it is expressed in terms of the two-particle correlation function $g(r)$ calculated for the ground state:

$$\langle 0 | \mathcal{A}_{\mathbf{q}} \mathcal{A}_{\mathbf{q}'}^\dagger | 0 \rangle = \frac{\nu}{N_\phi} \left[2\pi\nu \bar{g}(q) e^{q^2/2} + 1 \right] \delta_{\mathbf{q}', \mathbf{q}}. \quad (8)$$

Here $\bar{g}(q) = \frac{1}{(2\pi)^2} \int g(r) e^{-i\mathbf{q}\mathbf{r}} d^2r$ is the Fourier component. Function $g(r)$ is well known, e.g., in the case of Laughlin's state.^{12,13} If the ground state is presented in terms of the Hartree-Fock

model, we get the expression $2\pi\bar{g} = \left(N_\phi\delta_{\mathbf{q},0} - e^{-q^2/2}\right)$ which does not depend on ν . Besides, at odd-integer filling factors this Hartree-Fock expression becomes Fourier component of the *exact* correlation function. In the latter case one should also make the substitution $\nu \rightarrow \nu - 2n$ in Eq. (8), i.e. formally set $\nu = 1$ there.

3. The operator (7) obviously does not conserve the number of SEs. However, if the SWR is governed by this operator, the corresponding problem can not be solved in terms of a single-exciton study. Indeed, the SE interaction with the SRP incorporates the energy $U_{\text{x-SRP}} \sim ql_B\Delta/\Lambda$ (the SE possesses the dipole momentum $el_B[\mathbf{q} \times \hat{z}]$)². The SE momentum is estimated from the condition $\mathcal{E}_q \lesssim T$, and we therefore find that $U_{\text{x-SRP}} \ll \epsilon_Z, T$. Due to this inequality, the energy of annihilating exciton can not be transformed to anywhere. By analogy with Ref. 8, we study a coalescence process where initial double-exciton state $|i\rangle = \mathcal{Q}_{ab\mathbf{q}_1}^\dagger \mathcal{Q}_{ab\mathbf{q}_2}^\dagger |0\rangle$ transforms to final single-exciton state $|f\rangle = \mathcal{Q}_{ab\mathbf{q}'}^\dagger |0\rangle$ having the combined energy:

$$\epsilon_Z + \mathcal{E}_{q'} = 2\epsilon_Z + \mathcal{E}_{q_1} + \mathcal{E}_{q_2} \quad (9)$$

(c.f. also the Auger magnetoplasma relaxation considered in Ref. 14). At the same time, contrary to Ref. 8, there is no momentum conservation in this SWR channel. Thus the phase volume where the $X_{\mathbf{q}_1} + X_{\mathbf{q}_2} \rightarrow X_{\mathbf{q}'}$ transition is possible turns out to be much larger than that in the coalescence process of Ref. 8. This transition is governed by the Fermi golden rule probability: $w_{fi} = (2\pi/\hbar)|\mathcal{M}_{fi}|^2\delta(E_f - E_i)$, and our immediate task is to calculate the matrix element $\mathcal{M}_{fi} = \nu^{-3/2}\langle f|\hat{\varphi}|i\rangle$. (The factor $\nu^{-3/2}$ appears due to the normalization since norms of the $|i\rangle$ and $|f\rangle$ states are ν^2 and ν , respectively.)

We perform the calculation for relevant values of momenta $q_1, q_2, q' \ll 1$ which satisfy the conditions $\mathcal{E}_{q_1}, \mathcal{E}_{q_2} \lesssim T \lesssim 1$ K. (These inequalities correspond to $q_1, q_2, q' \ll 1/l_B$ in usual dimensional units). By employing exciton-operators' commutation rules⁷ and evident identities $\mathcal{Q}_{ab\mathbf{q}}|0\rangle \equiv \mathcal{B}_{\mathbf{q}}|0\rangle \equiv 0$ and $\langle 0|\mathcal{A}_{\mathbf{q}}|0\rangle \equiv \nu$, we obtain with the help of Eqs. (7)-(8) that

$$\mathcal{M}_{fi}(\mathbf{q}_1, \mathbf{q}_2, \mathbf{q}') = \frac{2\pi\nu^{1/2}}{N_\phi^{1/2}} \left[\sum_{j=1}^2 \bar{g}(\|\mathbf{q}_j - \mathbf{q}'\|) e^{(\mathbf{q}_j - \mathbf{q}')^2/2} \right] \sum_{\mathbf{q}} \bar{\varphi}(\mathbf{q})(iuq_+ - vq_-) \delta_{\mathbf{q}_1 + \mathbf{q}_2, \mathbf{q} + \mathbf{q}'}. \quad (10)$$

Besides, within our approximation, $\bar{g}(q)e^{q^2/2}$ should be replaced with $\bar{g}(q)e^{q^2/2}\Big|_{q \rightarrow 0}$. The latter quantity is equal to $-1/2\pi$ in the Hartree-Fock approach or $-1/2\pi\nu$ when calculated in the case of Laughlin's ground state describing the fractional QHF. So, for $\nu = 1, 1/3, 1/5, \dots$, replacing the terms in square brackets with $-1/\pi\nu$, we obtain a simple result:

$$|\mathcal{M}_{fi}(\mathbf{q}_1, \mathbf{q}_2, \mathbf{q}')|^2 = 4\pi \bar{K}(q) \frac{q^2(u^2 + v^2)}{\nu N_\phi^2} \Big|_{\mathbf{q} = \mathbf{q}_1 + \mathbf{q}_2 - \mathbf{q}'}. \quad (11)$$

It is used that the squared modulus of $\overline{\varphi}(\mathbf{q})$ may be expressed in terms of Fourier component of the correlator (4): $|\overline{\varphi}(\mathbf{q})|^2 = 2\pi K(q)/N_\phi$. In the Hartree-Fock model the expression (11) should be multiplied by ν^2 ; therefore the calculated relaxation rate would be by a factor of ν^2 slower. Notice also that if $\nu=3, 5, \dots$, one should formally set $\nu=1$ in Eqs. (10) and (11).

The SWR rate is defined as the difference between the fluxes of annihilating and created SEs. We assume that the thermodynamic equilibrium in the system of spin waves is established much faster than the spin-flip processes occur so that the rate is

$$R = \frac{1}{2} \sum_{\mathbf{q}_1, \mathbf{q}_2, \mathbf{q}'} \frac{2\pi}{\hbar} |\mathcal{M}_{fi}(\mathbf{q}_1, \mathbf{q}_2, \mathbf{q}')|^2 \delta(E_1 + E_2 - E') [n_1 n_2 (1 + n') - n' (1 + n_1) (1 + n_2)]. \quad (12)$$

The notations used here are $E_i = \epsilon_Z + \mathcal{E}_{q_i}$, $n_i = n(E_i)$ ($i = 1, 2$) and $E' = \epsilon_Z + \mathcal{E}_{q'}$, $n' = n(E')$, where the Bose distribution function is $n(E) = 1/(e^{(E-\mu)/T} - 1)$. The rate R is completely determined by Eqs. (11)-(12) and is a function of parameters B , T , and of the total number of SWs in the system: $N_x = \sum_{\mathbf{q}} n(\epsilon_Z + \mathcal{E}_q)$. In our case, when temperature is rather low, we can certainly use quadratic approximation for the “kinetic” exciton energy: $\mathcal{E}_q \approx q^2/2M_x$. Chemical potential μ is determined by the ratio of the exciton number and the total spin: $N_x(\mu) = \nu N_\phi/2 - S$. Calculating the quantity $N_x^{(0)} = N_x \Big|_{\mu=0}$, one obtains the equilibrium number of excitons. We will find the rate at the final stage of the relaxation process where $N_x - N_x^{(0)} \ll N_x^{(0)}$. So, by employing the quadratic approximation for the SE kinetic energy, and changing in Eqs. (11)-(12) from summations to integrations we obtain $R = (N_x - N_x^{(0)})/\tau_{\text{srp}}$, where

$$1/\tau_{\text{srp}} = \frac{(u^2 + v^2)M_x^3}{2\nu\pi\hbar} \left(\frac{\Delta\Lambda T}{l_B} \right)^2 (e^{-\epsilon_Z/T} - e^{-2\epsilon_Z/T}) F_{\text{SRP}}(\Lambda^2 M_x T/l_B^2, \epsilon_Z/T). \quad (13)$$

Here $F_{\text{SRP}}(\alpha, \beta)$ is a dimensionless function arising as a result of integrations over q_1 and q_2 and averaging over angles $\theta_1 = \mathbf{q}_1 \wedge \mathbf{q}'$ and $\theta_2 = \mathbf{q}_2 \wedge \mathbf{q}'$:

$$F_{\text{SRP}}(\alpha, \beta) = \int_0^\infty \int_0^\infty \frac{e^{-x-y} dx dy}{(1 - e^{-x-\beta})(1 - e^{-y-\beta})(1 - e^{-x-y-2\beta})} \\ \times \int_{-\pi}^\pi d\theta_1 \int_{-\pi}^\pi d\theta_2 r(x, y, \theta_1, \theta_2) \exp[-\alpha r(x, y, \theta_1, \theta_2)],$$

where $r(x, y, \theta_1, \theta_2) = x + y + \beta/2 - \sqrt{x+y+\beta}(\sqrt{x} \cos \theta_1 + \sqrt{y} \cos \theta_2) + \sqrt{xy} \cos(\theta_1 - \theta_2)$.

4. Now we calculate the numerical value of $1/\tau_{\text{srp}}$ at typical SRP parameters and compare it with inverse relaxation times $1/\tau_{e-e}$ and $1/\tau_{ph}$ governed by the inter-SEs' interaction mechanism⁸ and the SE-acoustic-phonon coupling.⁷ We carry out this analysis for the $\nu=1$ QHF assuming that $\Delta = 10$ K and $\Lambda = 40$ nm. The Zeeman splitting at $g = -0.44$ is $\epsilon_Z = 0.295B$ K (B is everywhere in Teslas), and the combination of SO parameters is estimated as $u^2 + v^2 = 10^{-3}/B$. The SE mass M_x might be calculated theoretically by using

general expressions for \mathcal{E}_q .^{2,3} Yet, the result depends on specific formfactor inherent in a given heterostructure due to finite thickness and it is therefore more convenient to extract M_x immediately from experiments. According to recent data available for currently used wide quantum wells,^{15,16} we estimate that $1/M_x = 9.24\sqrt{B}$ K. Using Eq. (13), we thus calculate $1/\tau_{\text{srp}}$ as a function of temperature T at given field B . The results are presented in Fig. 1 by dash curves. The dot and dash-dot curves correspond to the $1/\tau_{e-e}$ and $1/\tau_{ph}$ values given by formulas¹⁷

$$1/\tau_{e-e} = \frac{2}{\hbar}(u^2 + v^2)T (e^{-\epsilon_Z/T} - e^{-2\epsilon_Z/T}) F_{e-e}(\epsilon_Z/T), \quad (14)$$

where

$$F_{e-e}(\beta) = \iint_{xy > \beta^2/4} \frac{dxdy (x+y+\beta)e^{-x-y}}{(xy - \beta^2/4)^{1/2} (1 - e^{-\beta-x})(1 - e^{-\beta-y})(1 - e^{-2\beta-x-\beta})};$$

and

$$\tau_{ph}^{-1} = \frac{MT\epsilon_Z(u^2 + v^2)}{\hbar c_s p_0^3 l_B^2} \left[\frac{\gamma_1(\epsilon_Z/T)}{\tau_D} + 10 \frac{MT}{\tau_P} \left(\frac{\hbar c_s}{\epsilon_Z} \right)^4 \left(\frac{p_0}{l_B} \right)^2 \gamma_2(\epsilon_Z/T) \right], \quad (15)$$

where

$$\gamma_k(\beta) = (e^{2\beta} - e^\beta) \int_0^\infty \frac{e^x x^k dx}{(e^{\beta+x} - 1)^2}, \quad k=1, 2.$$

(See Ref. 7; the used material parameters characterizing the electron-phonon coupling are $c_s = 5.14 \cdot 10^5$ cm/s, $\tau_D = 0.8 \cdot 10^{-12}$ s⁻¹, $\tau_P = 35 \cdot 10^{-12}$ s⁻¹, and $p_0 = 2.52 \cdot 10^6$ cm⁻¹; both kinds of $e-ph$ interaction, deformation and polarization ones, are taken into account.)

It is seen from Fig. 1 that the SRP relaxation channel actually competes with other mechanisms in the experimentally relevant range of parameters: namely, at fields $B \leq 5$ and temperatures $T \sim 0.3 - 0.5$ K. We have indicated above that the basic advantage of the SRP channel, as compared to the $e-e$ one, consist in the absence of momentum conservation in the coalescence process. On the other hand, the SRP mechanisms is also determined by effective SE-SE collisions. Therefore the inverse relaxation time is proportional to the SE concentration and drops exponentially as $\sim \exp(-\epsilon_Z/T)$ with vanishing T [rather than as $\sim \exp(-2\epsilon_Z/T)$ which occurs for the $e-e$ mechanism due to the SEs momentum conservation!]. The phonon mechanism of SWR dominates at low temperatures due to its weak temperature dependence ($\sim T$), in spite of small value of the electron-phonon coupling constant in GaAs. The dependence on the filling factor in the case of integer QHF is only determined by the SE mass M_x because ν in Eq. (13) is formally set equal to unit. For fractional QHF there are both direct and indirect (through the mass M_x) dependences on ν .

Finally we calculate the combined inverse relaxation time determined by the SO interaction:

$$1/\tau_{\text{tot}} = 1/\tau_{\text{srp}} + 1/\tau_{e-e} + 1/\tau_{ph} \quad (16)$$

The result is presented by solid curves in Fig. 1. It is worth mentioning that it demonstrates a good agreement with the measured value $\tau_{\text{tot}} \simeq 10$ ns of Ref. 5 when calculated for parameters B and T corresponding to the experiment.

The authors acknowledge support of the RFBR and hospitality of the Max Planck Institute for Physics of Complex Systems (Dresden) where this work was partly carried out. The authors also thank S.V. Iordanskii and L.V. Kulik for discussion.

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 - ¹⁷ The corresponding formula for inverse relaxation time in Ref. 8 contains a misprint. Now we present the corrected result in Eq. (14).

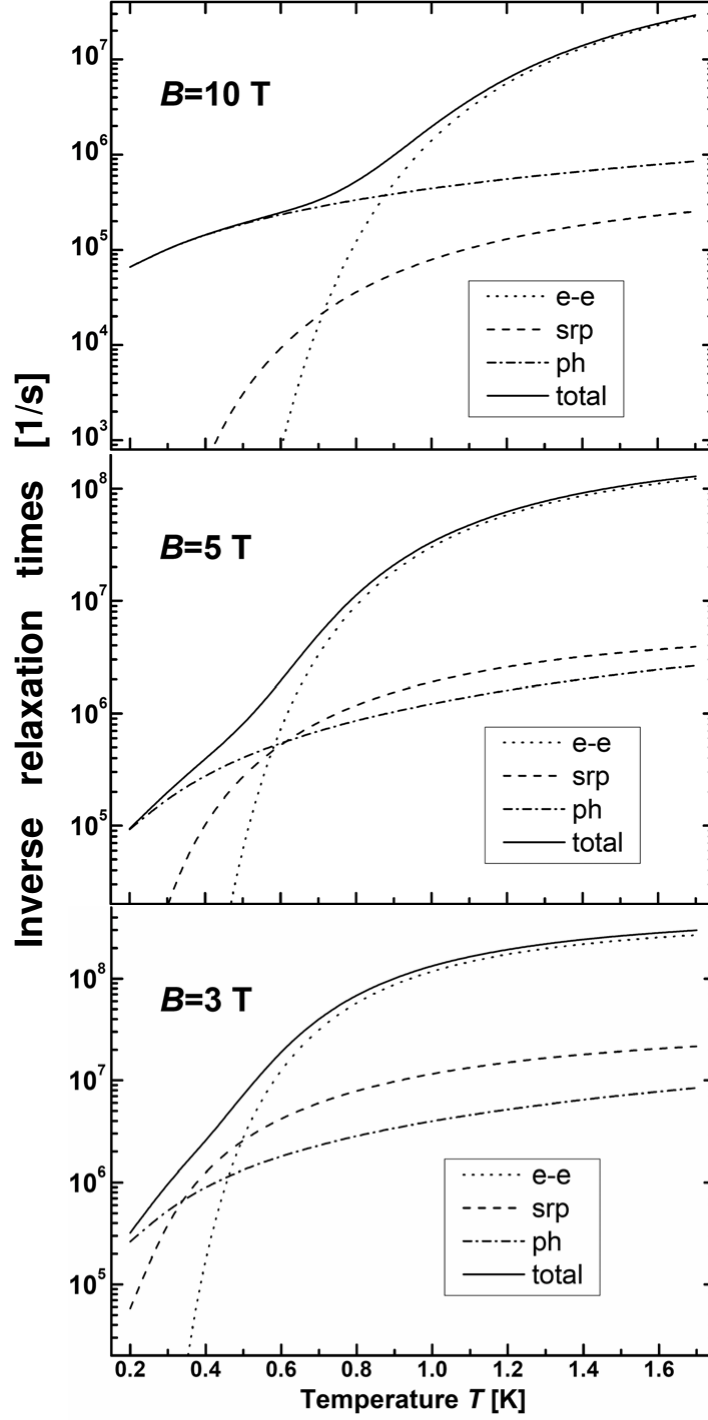


FIG. 1: Inverse SWR times against T calculated by using formulas (13)-(15) at $B = 3, 5, 10$ T. Specific material parameters are given in the text. Dash, dot, and dash-dot lines are for $1/\tau_{\text{srp}}$, $1/\tau_{\text{e-e}}$ and $1/\tau_{\text{ph}}$, respectively. Solid lines present the result of calculation of the combined inverse time (16).